```
=> b reg
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http://www.cas.org/support/stngen/stndoc/properties.html

=> d que sta l7
L1                STR
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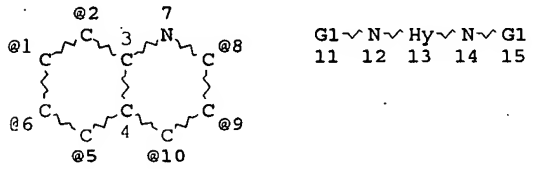


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VAR G1=8/9/10/5/6/1/2
NODE ATTRIBUTES:
DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED
ECOUNT  IS E4 C  E2 N  AT  13

GRAPH ATTRIBUTES:
RING(S) ARE ISOLATED OR EMBEDDED
NUMBER OF NODES IS  15

STEREO ATTRIBUTES: NONE
L7            44 SEA FILE=REGISTRY SSS FUL L1

100.0% PROCESSED  856075 ITERATIONS                          44 ANSWERS
SEARCH TIME: 00.00.06


=> b hcap
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FILE COVERS 1907 - 24 Jul 2007  VOL 147 ISS 5
FILE LAST UPDATED: 23 Jul 2007  (20070723/ED)
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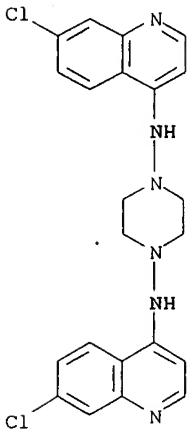
This file contains CAS Registry Numbers for easy and accurate substance identification.

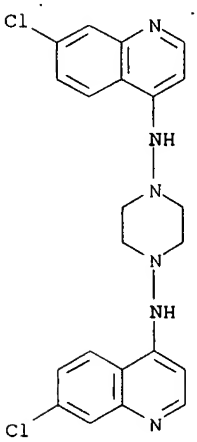
=> d bib abs hitstr l13 tot

L13 ANSWER 1 OF 2 HCAPLUS COPYRIGHT 2007 ACS on STN
AN 1971:435658 HCAPLUS
DN 75:35658
TI Antimalarials. "Distal" hydrazine derivatives of 7-chloroquinoline
AU Singh, Tara; Hoops, John F.; Biel, John H.; Hoya, Wallace K.; Stein, Robert George; Cruz, Deanna R.
CS Res. Lab., Aldrich Chem. Co., Inc., Milwaukee, WI, USA
SO Journal of Medicinal Chemistry (1971), 14(6), 532-5
CODEN: JMCMAR; ISSN: 0022-2623
DT Journal
LA English
GI For diagram(s), see printed CA Issue.
AB 7-Chloroquinolines (I) containing a hydrazine feature in the side chain attached at position 4, were prepared from 4,7-dichloroquinoline and 7-chloro-4-(3-bromo-1-methylpropylamino)quinoline by reacting with the required hydrazine, and were tested for the antimalarial activity against Plasmodium berghei in mice. 1,4-Bis(7-chloro-4-quinolylamino)-piperazine was the best, in which the end NH2 was substituted by a 2nd mol. of 7-chloroquinoline. It showed curative activity at 40 mg/kg, i.p., without toxicity even up to the maximum dose of 640 mg/kg. The I with a distal hydrazine, excluding active 1-[2-(7-chloro-4-quinolinylamino) - 2 - methylethyl] - 1 - methylhydrazine, were inactive, but were highly toxic. The I having a hydrazinium bromide feature, although found curative, were also quite toxic.
IT **23512-27-8P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 23512-27-8 HCAPLUS
CN 1,4-Piperazinediamine, N,N'-bis(7-chloro-4-quinolinyl)- (9CI) (CA INDEX NAME)



L13 ANSWER 2 OF 2 HCAPLUS COPYRIGHT 2007 ACS on STN
AN 1970:3335 HCAPLUS
DN 72:3335
TI Antimalarial substances. XVIII. Synthetic schistosomicides. 13. Antimalarial and antischistosomal effects of proximal hydrazine and hydroxylamine analogs of chloroquine and quinacrine
AU Elslager, Edward F.; Tendick, Frank H.; Werbel, Leslie M.; Worth, Donald F.
CS Med. and Sci. Affairs Div., Parke, Davis and Co., Ann Arbor, MI, USA
SO Journal of Medicinal Chemistry (1969), 12(5), 970-4
CODEN: JMCMAR; ISSN: 0022-2623
DT Journal

LA English
GI For diagram(s), see printed CA Issue.
AB Representative 4-(2,2-dialkylhydrazino)quinolines, 6 - chloro - 9 - (2,2 - dialkylhydrazino) - 2-methoxyacridines, 12-(2,2-dialkylhydrazino)benz[b]acridines, 2,-2'-(benz[c]acridin-7-ylhydrazono)diethanol,. 7-chloro-4 - [2- (dialkylamino)ethoxyamino]quinolines, and 6-chloro-9-[2- (dimethylamino)ethoxyamino]-2-methoxyacridine were synthesized to enable an assessment of the antiparasitic effects conferred by substituting a hydrazine or hydroxylamine moiety for the proximal amine function of chloroquine, quinacrine, and 7-[3-(octylamino)propylamino]benz[c]acridine relatives. The compds. were isolated in 3-92% yield by the condensation of 4,7-dichloroquinoline, 4-chloro-6-methoxyquinoline, 4-chloro-6-methoxyquinaldine, 6,9-dichloro-2-methoxyacridine, 12-chlorobenz[b]acridine, or 7-chlorobenz[c]acridine with the appropriate 1,1-dialkylhydrazine or 2-(dialkylamino)ethoxyamine in phenol or EtOH. Among them, 6-methoxy-4-(morpholinoamino)-quinaldine exhibited modest activity against Schistosoma mansoni in mice and effected a 28-51% reduction of live worms at drug-diet doses of 224-303 mg./kg. daily for 14 days. Six compds. were active against a normal strain of Plasmodium berghei in mice at doses ranging from 2.7-219 mg./kg./day for 6 days. 7-Chloro-4-(4-methyl-1-piperazinylamino)quinoline, and 4,4'-(1,4-piper-a zinediyldiimino)bis[7-chloroquinoline] (I) were approx. 28 and 27 times as potent as quinine, resp., against P. berghei, but I wa s highly cross-resistant with chloroquine. Structure-activity relations are discussed.
IT **23512-27-8P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 23512-27-8 HCAPLUS
CN 1,4-Piperazinediamine, N,N'-bis(7-chloro-4-quinolinyl)- (9CI) (CA INDEX NAME)



=> d his

(FILE 'HCAPLUS' ENTERED AT 11:18:23 ON 24 JUL 2007)
DEL HIS Y

FILE 'REGISTRY' ENTERED AT 11:36:22 ON 24 JUL 2007
L1 STR
L2 0 L1

FILE 'HCAPLUS' ENTERED AT 11:39:10 ON 24 JUL 2007
L3 1 (US20040053966 OR US6645964)/PN OR (US2003-658394 OR FR2000-105

FILE 'REGISTRY' ENTERED AT 11:42:58 ON 24 JUL 2007

FILE 'HCAPLUS' ENTERED AT 11:42:58 ON 24 JUL 2007
L4 TRA L3 1- RN : 73 TERMS

FILE 'REGISTRY' ENTERED AT 11:42:58 ON 24 JUL 2007

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L5            73 SEA L4
L6             2 L5 AND (NC2NC2 OR N2C4 OR NCNC3)/ES
L7            44 L1 FULL
                 SAV TEM J394C1/A L7

     FILE 'HCAPLUS' ENTERED AT 12:58:21 ON 24 JUL 2007
L8            11 L7
L9             1 L8 AND L3
L10           10 L8 NOT L9
                 SEL HIT RN L10 6-10

     FILE 'REGISTRY' ENTERED AT 13:00:13 ON 24 JUL 2007
L11            5 E27-31

     FILE 'STNGUIDE' ENTERED AT 13:01:14 ON 24 JUL 2007

     FILE 'REGISTRY' ENTERED AT 13:03:34 ON 24 JUL 2007
L12            1 L11 AND C22H20CL2N6

     FILE 'HCAPLUS' ENTERED AT 13:04:39 ON 24 JUL 2007
L13            2 L12

     FILE 'HCAOLD' ENTERED AT 13:05:20 ON 24 JUL 2007
L14            0 L7


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